

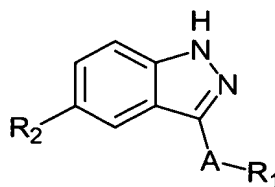
Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application.

Listing of Claims

1-5. (Canceled)

6. (Previously presented) A compound having the structure:



or a pharmaceutically acceptable salt thereof, wherein:

A is $-(CH_2)_bC \equiv C(CH_2)_c-$;

R₁ is aryl, heteroaryl or heterocycle fused to phenyl, each being optionally substituted with one to four substituents independently selected from R₃;

R₂ is -R₃, -R₄, $-(CH_2)_bC(=O)R_5$, $-(CH_2)_bC(=O)OR_5$, $-(CH_2)_bC(=O)NR_5R_6$, $-(CH_2)_bC(=O)NR_5(CH_2)_cC(=O)R_6$, $-(CH_2)_bNR_5C(=O)R_6$, $-(CH_2)_bNR_5C(=O)NR_6R_7$, $-(CH_2)_bNR_5R_6$, $-(CH_2)_bOR_5$, $-(CH_2)_bSO_dR_5$ or $-(CH_2)_bSO_2NR_5R_6$;

b and *c* are the same or different and at each occurrence independently selected from 0, 1, 2, 3 or 4;

d is at each occurrence 0, 1 or 2;

R₃ is at each occurrence independently halogen, hydroxy, carboxy, alkyl, alkoxy, haloalkyl, acyloxy, thioalkyl, sulfinylalkyl, sulfonylalkyl, hydroxyalkyl, aryl, substituted aryl, arylalkyl, substituted arylalkyl, heterocycle, substituted heterocycle, heterocyclealkyl, substituted heterocyclealkyl, $-C(=O)OR_8$, $-C(=O)R_8$, $-C(O)NR_8R_9$, $-C(=O)NR_8OR_9$, $-SO_2NR_8R_9$, $-NR_8SO_2R_9$, $-CN$, $-NO_2$, $-NR_8R_9$, $-NR_8C(=O)R_9$, $-NR_8C(=O)(CH_2)_bOR_9$, $-NR_8C(=O)(CH_2)_bR_9$, $-O(CH_2)_bNR_8R_9$, or heterocycle fused to phenyl;

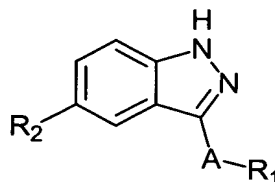
R₄ is alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, each being optionally substituted with one to four substituents independently selected from R₃, or R₄ is halogen or hydroxy;

R₅, R₆ and R₇ are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, wherein each of R₅, R₆ and R₇ are optionally substituted with one to four substituents independently selected from R₃; and

R₈ and R₉ are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle, or heterocyclealkyl, or R₈ and R₉ taken together with the atom or atoms to which they are bonded form a heterocycle, wherein each of R₈, R₉, and R₈ and R₉ taken together to form a heterocycle are optionally substituted with one to four substituents independently selected from R₃.

7-9. (Canceled)

10. (Previously presented) A compound having the structure:



or a pharmaceutically acceptable salt thereof, wherein:

A is a direct bond, $-(CH_2)_a-$, $-(CH_2)_bCH=CH(CH_2)_c-$, or $-(CH_2)_bC \equiv C(CH_2)_c-$;

R₁ is aryl, heteroaryl or heterocycle fused to phenyl, each being optionally substituted with one to four substituents independently selected from R₃;

R₂ is $-(CH_2)_bC(=O)R_5$;

a is 1, 2, 3, 4, 5 or 6;

b and c are the same or different and at each occurrence independently selected from 0, 1, 2, 3 or 4;

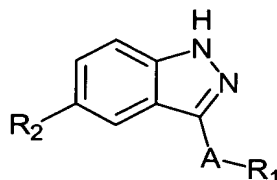
R₃ is at each occurrence independently halogen, hydroxy, carboxy, alkyl, alkoxy, haloalkyl, acyloxy, thioalkyl, sulfinylalkyl, sulfonylalkyl, hydroxyalkyl, aryl, substituted aryl, arylalkyl, substituted arylalkyl,

heterocycle, substituted heterocycle, heterocyclealkyl, substituted heterocyclealkyl, $-C(=O)OR_8$, $-C(=O)R_8$, $-C(O)NR_8R_9$, $-C(=O)NR_8OR_9$, $-SO_2NR_8R_9$, $-NR_8SO_2R_9$, $-CN$, $-NO_2$, $-NR_8R_9$, $-NR_8C(=O)R_9$, $-NR_8C(=O)(CH_2)_bOR_9$, $-NR_8C(=O)(CH_2)_bR_9$, $-O(CH_2)_bNR_8R_9$, or heterocycle fused to phenyl;

R_5 , R_6 and R_7 are the same or different and at each occurrence independently alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, wherein each of R_5 , R_6 and R_7 are optionally substituted with one to four substituents independently selected from R_3 ; and

R_8 and R_9 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle, or heterocyclealkyl, or R_8 and R_9 taken together with the atom or atoms to which they are bonded form a heterocycle, wherein each of R_8 , R_9 , and R_8 and R_9 taken together to form a heterocycle are optionally substituted with one to four substituents independently selected from R_3 .

11. (Previously presented) A compound having the structure:



or a pharmaceutically acceptable salt thereof, wherein:

A is a direct bond, $-(CH_2)_a-$, $-(CH_2)_bCH=CH(CH_2)_c-$, or $-(CH_2)_bC \equiv C(CH_2)_c-$;

R_1 is aryl, heteroaryl or heterocycle fused to phenyl, each being optionally substituted with one to four substituents independently selected from R_3 ;

R_2 is $-(CH_2)_bC(=O)NR_5R_6$;

a is 1, 2, 3, 4, 5 or 6;

b and c are the same or different and at each occurrence independently selected from 0, 1, 2, 3 or 4;

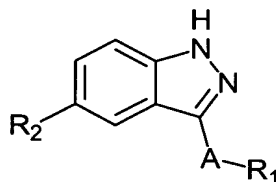
R_3 is at each occurrence independently halogen, hydroxy, carboxy, alkyl, alkoxy, haloalkyl, acyloxy, thioalkyl, sulfinylalkyl, sulfonylalkyl, hydroxyalkyl, aryl, substituted aryl, arylalkyl, substituted arylalkyl,

heterocycle, substituted heterocycle, heterocyclealkyl, substituted heterocyclealkyl, $-C(=O)OR_8$, $-C(=O)R_8$, $-C(O)NR_8R_9$, $-C(=O)NR_8OR_9$, $-SO_2NR_8R_9$, $-NR_8SO_2R_9$, $-CN$, $-NO_2$, $-NR_8R_9$, $-NR_8C(=O)R_9$, $-NR_8C(=O)(CH_2)_bOR_9$, $-NR_8C(=O)(CH_2)_bR_9$, $-O(CH_2)_bNR_8R_9$, or heterocycle fused to phenyl;

R_5 , R_6 and R_7 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, wherein each of R_5 , R_6 and R_7 are optionally substituted with one to four substituents independently selected from R_3 ; and

R_8 and R_9 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle, or heterocyclealkyl, or R_8 and R_9 taken together with the atom or atoms to which they are bonded form a heterocycle, wherein each of R_8 , R_9 , and R_8 and R_9 taken together to form a heterocycle are optionally substituted with one to four substituents independently selected from R_3 .

12. (Previously presented) A compound having the structure:



or a pharmaceutically acceptable salt thereof, wherein:

A is a direct bond, $-(CH_2)_a-$, $-(CH_2)_bCH=CH(CH_2)_c-$, or $-(CH_2)_bC\equiv C(CH_2)_c-$;

R_1 is aryl, heteroaryl or heterocycle fused to phenyl, each being optionally substituted with one to four substituents independently selected from R_3 ;

R_2 is $-(CH_2)_bNR_5C(=O)R_6$;

a is 1, 2, 3, 4, 5 or 6;

b and c are the same or different and at each occurrence independently selected from 0, 1, 2, 3 or 4;

R_3 is at each occurrence independently halogen, hydroxy, carboxy, alkyl, alkoxy, haloalkyl, acyloxy, thioalkyl, sulfinylalkyl, sulfonylalkyl, hydroxyalkyl, aryl, substituted aryl, arylalkyl, substituted arylalkyl,

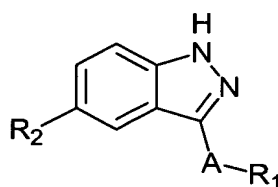
heterocycle, substituted heterocycle, heterocyclealkyl, substituted heterocyclealkyl, $-\text{C}(=\text{O})\text{OR}_8$, $-\text{C}(=\text{O})\text{R}_8$, $-\text{C}(\text{O})\text{NR}_8\text{R}_9$, $-\text{C}(=\text{O})\text{NR}_8\text{OR}_9$, $-\text{SO}_2\text{NR}_8\text{R}_9$, $-\text{NR}_8\text{SO}_2\text{R}_9$, $-\text{CN}$, $-\text{NO}_2$, $-\text{NR}_8\text{R}_9$, $-\text{NR}_8\text{C}(=\text{O})\text{R}_9$, $-\text{NR}_8\text{C}(=\text{O})(\text{CH}_2)_b\text{OR}_9$, $-\text{NR}_8\text{C}(=\text{O})(\text{CH}_2)_b\text{R}_9$, $-\text{O}(\text{CH}_2)_b\text{NR}_8\text{R}_9$, or heterocycle fused to phenyl;

R_5 , R_6 and R_7 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, wherein each of R_5 , R_6 and R_7 are optionally substituted with one to four substituents independently selected from R_3 ; and

R_8 and R_9 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle, or heterocyclealkyl, or R_8 and R_9 taken together with the atom or atoms to which they are bonded form a heterocycle, wherein each of R_8 , R_9 , and R_8 and R_9 taken together to form a heterocycle are optionally substituted with one to four substituents independently selected from R_3 .

13-17. (Canceled)

18. (Previously presented) A compound having the structure:



or a pharmaceutically acceptable salt thereof, wherein:

A is a direct bond, $-(\text{CH}_2)_a-$, $-(\text{CH}_2)_b\text{CH}=\text{CH}(\text{CH}_2)_c-$, or $-(\text{CH}_2)_b\text{C}\equiv\text{C}(\text{CH}_2)_c-$;

R_1 is aryl, heteroaryl or heterocycle fused to phenyl, each being optionally substituted with one to four substituents independently selected from R_3 ;

R_2 is R_4 ;

a is 1, 2, 3, 4, 5 or 6;

b and c are the same or different and at each occurrence independently selected from 0, 1, 2, 3 or 4;

R₃ is at each occurrence independently halogen, hydroxy, carboxy, alkyl, alkoxy, haloalkyl, acyloxy, thioalkyl, sulfinylalkyl, sulfonylalkyl, hydroxyalkyl, aryl, substituted aryl, arylalkyl, substituted arylalkyl, heterocycle, substituted heterocycle, heterocyclealkyl, substituted heterocyclealkyl, -C(=O)OR₈, -C(=O)R₈, -C(O)NR₈R₉, -C(=O)NR₈OR₉, -SO₂NR₈R₉, -NR₈SO₂R₉, -CN, -NO₂, -NR₈R₉, -NR₈C(=O)R₉, -NR₈C(=O)(CH₂)_bOR₉, -NR₈C(=O)(CH₂)_bR₉, -O(CH₂)_bNR₈R₉, or heterocycle fused to phenyl;

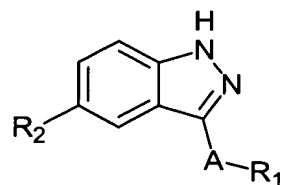
R₄ is 3-triazolyl, optionally substituted at its 5-position with:

- (a) a C₁-C₄ straight or branched chain alkyl group optionally substituted with a hydroxyl, methylamino, dimethylamino or 1-pyrrolidinyl group; or
- (b) a 2-pyrrolidinyl group;

R₅, R₆ and R₇ are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, wherein each of R₅, R₆ and R₇ are optionally substituted with one to four substituents independently selected from R₃; and

R₈ and R₉ are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle, or heterocyclealkyl, or R₈ and R₉ taken together with the atom or atoms to which they are bonded form a heterocycle, wherein each of R₈, R₉, and R₈ and R₉ taken together to form a heterocycle are optionally substituted with one to four substituents independently selected from R₃.

19. (Previously presented) A compound having the structure:



or a pharmaceutically acceptable salt thereof, wherein:

A is a direct bond, -(CH₂)_a-, -(CH₂)_bCH=CH(CH₂)_c-, or -(CH₂)_bC≡C(CH₂)_c-;

R_1 is aryl, heteroaryl or heterocycle fused to phenyl, each being optionally substituted with one to four substituents independently selected from R_3 ;

R_2 is R_4 ;

a is 1, 2, 3, 4, 5 or 6;

b and c are the same or different and at each occurrence independently selected from 0, 1, 2, 3 or 4;

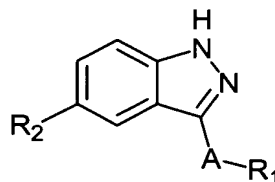
R_3 is at each occurrence independently halogen, hydroxy, carboxy, alkyl, alkoxy, haloalkyl, acyloxy, thioalkyl, sulfinylalkyl, sulfonylalkyl, hydroxyalkyl, aryl, substituted aryl, arylalkyl, substituted arylalkyl, heterocycle, substituted heterocycle, heterocyclealkyl, substituted heterocyclealkyl, $-C(=O)OR_8$, $-C(=O)R_8$, $-C(O)NR_8R_9$, $-C(=O)NR_8OR_9$, $-SO_2NR_8R_9$, $-NR_8SO_2R_9$, $-CN$, $-NO_2$, $-NR_8R_9$, $-NR_8C(=O)R_9$, $-NR_8C(=O)(CH_2)_bOR_9$, $-NR_8C(=O)(CH_2)_bR_9$, $-O(CH_2)_bNR_8R_9$, or heterocycle fused to phenyl;

R_4 is tetrazole;

R_5 , R_6 and R_7 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, wherein each of R_5 , R_6 and R_7 are optionally substituted with one to four substituents independently selected from R_3 ; and

R_8 and R_9 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle, or heterocyclealkyl, or R_8 and R_9 taken together with the atom or atoms to which they are bonded form a heterocycle, wherein each of R_8 , R_9 , and R_8 and R_9 taken together to form a heterocycle are optionally substituted with one to four substituents independently selected from R_3 .

20. (Previously presented) A compound having the structure:



or a pharmaceutically acceptable salt thereof, wherein:

A is a direct bond, $-(CH_2)_a-$, $-(CH_2)_bCH=CH(CH_2)_c-$, or $-(CH_2)_bC\equiv C(CH_2)_c-$;

R_1 is aryl, heteroaryl or heterocycle fused to phenyl, each being optionally substituted with one to four substituents independently selected from R_3 ;

R_2 is R_4 ;

a is 1, 2, 3, 4, 5 or 6;

b and c are the same or different and at each occurrence independently selected from 0, 1, 2, 3 or 4;

R_3 is at each occurrence independently halogen, hydroxy, carboxy, alkyl, alkoxy, haloalkyl, acyloxy, thioalkyl, sulfinylalkyl, sulfonylalkyl, hydroxyalkyl, aryl, substituted aryl, arylalkyl, substituted arylalkyl, heterocycle, substituted heterocycle, heterocyclealkyl, substituted heterocyclealkyl, $-C(=O)OR_8$, $-C(=O)R_8$, $-C(O)NR_8R_9$, $-C(=O)NR_8OR_9$, $-SO_2NR_8R_9$, $-NR_8SO_2R_9$, $-CN$, $-NO_2$, $-NR_8R_9$, $-NR_8C(=O)R_9$, $-NR_8C(=O)(CH_2)_bOR_9$, $-NR_8C(=O)(CH_2)_bR_9$, $-O(CH_2)_bNR_8R_9$, or heterocycle fused to phenyl;

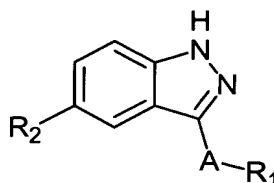
R_4 is imidazole;

R_5 , R_6 and R_7 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, wherein each of R_5 , R_6 and R_7 are optionally substituted with one to four substituents independently selected from R_3 ; and

R_8 and R_9 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle, or heterocyclealkyl, or R_8 and R_9 taken together with the atom or atoms to which they are bonded form a heterocycle, wherein each of R_8 , R_9 , and R_8 and R_9 taken together to form a heterocycle are optionally substituted with one to four substituents independently selected from R_3 .

21-73. (Canceled)

74. (Previously presented) A compound having the structure:



or a pharmaceutically acceptable salt thereof, wherein:

-A-R₁ is phenyl, optionally substituted with one to four substituents independently selected from halogen, alkoxy, -NR₈C(=O)R₉, -C(=O)NR₈R₉, and -O(CH₂)_bNR₈R₉;

R₂ is 3-triazolyl or 5-tetrazolyl;

a is 1, 2, 3, 4, 5 or 6;

b is 2 or 3;

c is at each occurrence 0, 1, 2, 3 or 4;

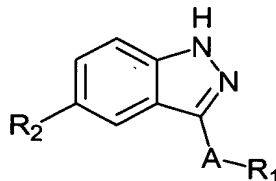
R₃ is at each occurrence independently halogen, hydroxy, carboxy, alkyl, alkoxy, haloalkyl, acyloxy, thioalkyl, sulfinylalkyl, sulfonylalkyl, hydroxyalkyl, aryl, substituted aryl, arylalkyl, substituted arylalkyl, heterocycle, substituted heterocycle, heterocyclealkyl, substituted heterocyclealkyl, -C(=O)OR₈, -C(=O)R₈, -C(O)NR₈R₉, -C(=O)NR₈OR₉, -SO₂NR₈R₉, -NR₈SO₂R₉, -CN, -NO₂, -NR₈R₉, -NR₈C(=O)R₉, -NR₈C(=O)(CH₂)_bOR₉, -NR₈C(=O)(CH₂)_bR₉, -O(CH₂)_bNR₈R₉, or heterocycle fused to phenyl;

R₅, R₆ and R₇ are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, wherein each of R₅, R₆ and R₇ are optionally substituted with one to four substituents independently selected from R₃; and

R₈ and R₉ are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle, or heterocyclealkyl, or R₈ and R₉ taken together with the atom or atoms to which they are bonded form a heterocycle, wherein each of R₈, R₉, and R₈ and R₉ taken together to form a heterocycle are optionally substituted with one to four substituents independently selected from R₃.

75-84. (Canceled)

85. (Previously presented) A compound having the structure:



or a pharmaceutically acceptable salt thereof, wherein:

A is a direct bond, $-(CH_2)_a-$, $-(CH_2)_bCH=CH(CH_2)_c-$, or $-(CH_2)_bC\equiv C(CH_2)_c-$;

R_1 is aryl, heteroaryl or heterocycle fused to phenyl, each being optionally substituted with one to four substituents independently selected from R_3 ;

R_2 is R_4 ;

a is 1, 2, 3, 4, 5 or 6;

b and c are the same or different and at each occurrence independently selected from 0, 1, 2, 3 or 4;

R_3 is at each occurrence independently halogen, hydroxy, carboxy, alkyl, alkoxy, haloalkyl, acyloxy, thioalkyl, sulfinylalkyl, sulfonylalkyl, hydroxyalkyl, aryl, substituted aryl, arylalkyl, substituted arylalkyl, heterocycle, substituted heterocycle, heterocyclealkyl, substituted heterocyclealkyl, $-C(=O)OR_8$, $-C(=O)R_8$, $-C(O)NR_8R_9$, $-C(=O)NR_8OR_9$, $-SO_2NR_8R_9$, $-NR_8SO_2R_9$, $-CN$, $-NO_2$, $-NR_8R_9$, $-NR_8C(=O)R_9$, $-NR_8C(=O)(CH_2)_bOR_9$, $-NR_8C(=O)(CH_2)_bR_9$, $-O(CH_2)_bNR_8R_9$, or heterocycle fused to phenyl;

R_4 is 3-triazolyl, optionally substituted at its 5-position with:

(a) methyl, n-propyl, isopropyl, 1-hydroxyethyl, 3-hydroxypropyl, methylaminomethyl, dimethylaminomethyl, 1-(dimethylamino)ethyl, 1-pyrrolidinylmethyl or 2-pyrrolidinyl;

R_5 , R_6 and R_7 are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle or heterocyclealkyl, wherein each of R_5 , R_6 and R_7 are optionally substituted with one to four substituents independently selected from R_3 ; and

R₈ and R₉ are the same or different and at each occurrence independently hydrogen, alkyl, aryl, arylalkyl, heterocycle, or heterocyclealkyl, or R₈ and R₉ taken together with the atom or atoms to which they are bonded form a heterocycle, wherein each of R₈, R₉, and R₈ and R₉ taken together to form a heterocycle are optionally substituted with one to four substituents independently selected from R₃.

86-88. (Canceled)

89. (Previously presented) A composition comprising the compound of claim 6 and a pharmaceutically acceptable carrier.

90. (Previously presented) A composition comprising the compound of claim 10 and a pharmaceutically acceptable carrier.

91. (Previously presented) A composition comprising the compound of claim 11 and a pharmaceutically acceptable carrier.

92. (Previously presented) A composition comprising the compound of claim 12 and a pharmaceutically acceptable carrier.

93-97. (Canceled)

98. (Previously presented) A composition comprising the compound of claim 18 and a pharmaceutically acceptable carrier.

99. (Previously presented) A composition comprising the compound of claim 19 and a pharmaceutically acceptable carrier.

100. (Previously presented) A composition comprising the compound of claim 20 and a pharmaceutically acceptable carrier.

101-103. (Canceled)

104. (Previously presented) A composition comprising the compound of claim 74 and a pharmaceutically acceptable carrier.

105. ((Previously presented) A composition comprising the compound of claim 85 and a pharmaceutically acceptable carrier.

106. (Previously presented) A compound of claim 6, wherein the compound is:

3-(2-phenylethynyl)-1H-indazole-5-carboxamide, or a pharmaceutically acceptable salt thereof.

107. (Previously presented) A compound of claim 10, wherein the compound is:

1-{{(3-(4-fluorophenyl)-1H-indazol-5-yl)carbonyl} piperidine-4-carboxylic acid;
3-(4-fluorophenyl)(1H-indazol-5-yl) pyrrolidinyl ketone;
3-(4-fluorophenyl)(1H-indazol-5-yl)piperazinyl ketone;
1-(3-(4-fluorophenyl)-1H-indazol-5-yl)-2-phenylethan-1-one;
1-(3-(4-fluorophenyl)-1H-indazol-5-yl)ethan-1-one; or a pharmaceutically acceptable salt thereof.

108. (Previously presented) A compound of claim 11, wherein the compound is:

3-(4-fluorophenyl)-1H-indazole-5-carboxamide;
(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-benzamide;
N-(2-(dimethylamino)ethyl)3-(4-fluorophenyl) (1H-indazol-5-yl)carboxamide;
methyl 4-{{(3-(4-fluorophenyl)-1H-indazol-5-yl)carbonylamino}benzoate;
4-{{(3-(4-fluorophenyl)-1H-indazol-5-yl)carbonylamino}benzoic acid;
4-{{(3-(4-fluorophenyl)-1H-indazole-5-yl)carbonylamino}benzamide;
(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(2-pyridyl)carboxamide;
(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(3-pyridyl)carboxamide;
(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(4-pyridyl)carboxamide;
tert-butyl 3-{{(3-(4-fluorophenyl)-1H-indazol-5-yl)carbonylamino}propanoate;
(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(3-hydroxyphenyl)carboxamide;
3-{{(3-(4-fluorophenyl)-1H-Indazol-5-yl)carbonylamino}propanoic acid;
(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(3-nitrophenyl)carboxamide;
tert-butyl-2-{{(3-(4-fluorophenyl)-1H-indazol-5-yl)carbonylamino}acetate;
4-{{(3-(4-fluorophenyl)-1H-indazol-5-yl)carbonylamino} butanoic acid;

N-(3-aminophenyl)(3-(4-fluorophenyl)(1H-indazol-5-yl))carboxamide;
 2-((3-(4-fluorophenyl)-1H-indazol-5-yl)carbonylamino}acetic acid;
 5-((3-(4-fluorophenyl)-1H-indazol-5-yl)carbonylamino}pentanoic acid;
 4-((3-(4-fluorophenyl)-1H-indazol-5-yl)carbonylamino}methyl)benzoic acid;
 (3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(4-pyridylmethyl)carboxamide;
 2-(4-((3-(4-fluorophenyl)-1H-indazol-5-yl)carbonylamino}phenyl)acetic acid;
 (3-(4-fluorophenyl)(1H-indazol-5-yl))-N,N-dimethylcarboxamide;
 (3-(4-fluorophenyl)(1H-indazol-5-yl))-N-methylcarboxamide;
 N-(3-aminoethyl)(3-(4-fluorophenyl)(1H-indazol-5-yl))carboxamide;
 N-(3-aminopropyl)(3-(4-fluorophenyl)(1H-indazol-5-yl))carboxamide;
 (3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(2-hydroxypropyl)carboxamide;
 N-(2H-1,2,3,4-tetrazol-5-yl)(3-(4-fluorophenyl)(1H-indazol-5-yl))carboxamide;
 {3-(4-fluorophenyl)(1H-indazol-5-yl)}-N-(3-morpholin-4-ylpropyl)carboxamide;
 (3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(3-pyridylmethyl)carboxamide;
 N-(((2R)-2-hydroxycyclohexyl)methyl)(3-(4-fluorophenyl)(1H-indazole-5-yl))carboxamide;
 (3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(2-(1-methylimidazol-5-yl)ethyl)carboxamide);
 (3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(2-pyridylmethyl)carboxamide;
 N-(2-carbamoyl)ethyl(3-(4-fluorophenyl)(1H-indazol-5-yl))carboxamide;
 N-(3-carbamoylpropyl)(3-(4-fluorophenyl)(1H-indazol-5-yl))carboxamide;
 3-(4-methoxyphenyl)-1H-indazole-5-carboxamide;
 3-(4-hydroxyphenyl)-1H-indazole-5-carboxamide;
 3-(2-naphthyl)-1H-indazole-5-carboxamide;
 3-benzo(b)thiophen-2-yl-1H-indazole-5-carboxamide;
 3-benzo(d)furan-2-yl-1H-indazole-5-carboxamide;
 3-(3-(methylethyl)phenyl)-1H-indazole-5-carboxamide;
 3-(4-(dimethylamino)phenyl)-1H-indazole-5-carboxamide;
 3-(3-furyl)-1H-indazole-5-carboxamide;
 3-{4-(2-(dimethylamino)ethoxy)phenyl}-1H-indazole-5-carboxamide;
 3-(3,4-dimethoxyphenyl)-1H-indazole-5-carboxamide;
 3-(3-aminophenyl)-1H-indazole-5-carboxamide;
 3-(2H-benzo(d)1,3-dioxolen-5-yl)-1H-indazole-5-carboxamide;

(3-benzo(d)furan-2-yl(1H-indazol-5-yl))-N-(methylethyl)carboxamide;
 (3-benzo(d)furan-2-yl(1H-indazol-5-yl))-N-(2-methoxyethyl)carboxamide;
 (3-benzo(d)furan-2-yl(1H-indazol-5-yl))-N-(2-(dimethylamino)ethyl)carboxamide;
 (3-benzo(d)furan-2-yl(1H-indazol-5-yl))-N-(4-(dimethylamino)butyl)carboxamide;
 (3-benzo(d)furan-2-yl(1H-indazol-5-yl))-N-(3-(dimethylamino)propyl)carboxamide;
 (3-benzo(d)furan-2-yl(1H-indazol-5-yl))-N-(2-methylpropyl)carboxamide;
 (3-benzo(d)furan-2-yl(1H-indazol-5-yl))-N-methylcarboxamide;
 3-(3-(3-pyridylcarbonylamino)phenyl)-1H-indazole-5-carboxamide;
 3-(3-(2-methoxyacetylaminophenyl)-1H-indazole-5-carboxamide;
 3-(3-(4-piperidylcarboxyamino)phenyl)-1H-indazole-5-carboxamide;
 (1S)-1-{N-(3-(5-carbamoyl(1H-indazol-3-yl))phenyl)carbamoyl}ethyl acetate;
 3-{3-(2-methoxyethyl)amino}phenyl}-1H-indazole-5-carboxamide;
 3-(3-(3-piperidylpropanoylamino)phenyl)-1H-indazole-5-carboxamide;
 3-(3-(2-furylecarbonylamino)phenyl)-1H-indazole-5-carboxamide;
 3-{3-(2-(dimethylamino)acetylaminophenyl)-1H-indazole-5-carboxamide;
 3-(3-(2-phenylacetylaminophenyl)-1H-Indazole-5-carboxamide;
 3-{3-(2-(4-methoxyphenyl)acetylaminophenyl)-1H-indazole-5-carboxamide;
 3-{3-(2-(2-methyl-1,3-thiazol-5-yl)acetylaminophenyl)-1H-indazole-5-
 carboxamide;
 3-(3-(oxolan-3-yl-carbonylamino)phenyl)-1H-indazole-5-carboxamide;
 3-(3-(2-(3-thienyl)acetylaminophenyl)-1H-indazole-5-carboxamide;
 3-(3-(2-thienylcarbonylamino)phenyl)-1H-indazole-5-carboxamide;
 3-(3-(2-(4-pyridyl)acetylaminophenyl)-1H-Indazole-5-carboxamide;
 3-(3-(2-(2-pyridyl)acetylaminophenyl)-1H-Indazole-5-carboxamide;
 3-{3-(2-(4-fluorophenyl)acetylaminophenyl)-1H-indazole-5-carboxamide;
 3-(3-(cyclopropylcarbonylamino)phenyl)-1H-indazole-5-carboxamide;
 3-{3-((3-hydroxyphenyl)carbonylamino)phenyl}-1H-indazole-5-carboxamide;
 3-{3-(2-(2,4-dichlorophenyl)acetylaminophenyl)-1H-indazole-5-carboxamide;
 3-(3-{2-(4-(trifluoromethyl)phenyl)acetylaminophenyl)-1H-indazole-5-
 carboxamide;
 3-(3-{2-(4-(dimethylamino)phenyl)acetylaminophenyl)-1H-indazole-5-
 carboxamide;

3-{3-(2-(2-chloro-4-fluorophenyl) acetylamino)phenyl}-1H-indazole-5-carboxamide;

3-{3-(2-(4-chlorophenyl)acetylamino)phenyl}-1H-indazole-5-carboxamide;

3-(3-(3-phenylpropanoylamino)phenyl)-1H-indazole-5-carboxamide;

3-{3-(3-(4-fluorophenyl)propanoylamino)phenyl}-1H-indazole-5-carboxamide;

3-{3-(2-(3,4-difluorophenyl)acetylamino)phenyl}-1H-indazole-5-carboxamide;

3-{3-(2-(2-fluorophenyl) acetylamino)phenyl}-1H-indazole-5-carboxamide;

3-(3-(2-phenylpropanoylamino)phenyl)-1H-indazole-5-carboxamide;

3-(3-(2-piperidylethoxy)phenyl)-1H-indazole-5-carboxamide;

N-ethyl-3-((3-(4-fluorophenyl)(1H-indazol-5-yl))carbonylamino) propanamide;

(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(3-methoxypropyl)carboxamide;

3-{3-(N-(2-piperidylethyl)carbamoyl)phenyl}-1H-indazole-5-carboxamide;

(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(2-hydroxyethyl)carboxamide;

(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(3-hydroxypropyl)carboxamide;

(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(2-methoxyethyl)carboxamide;

(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(oxolan-2-ylmethyl)carboxamide;

3-(2H, 3H-benzo(e)1,4-dioxin-6-yl)-1H-indazole-5-carboxamide;

3-(3-quinolyl)-1H-indazole-5-carboxamide;

3-(6-methoxy-2-naphthyl)-1H-indazole-5-carboxamide;

3-(2,3-dihydrobenzo(b)furan-5-yl)-1H-indazole-5-carboxamide;

(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(3-oxo-3-pyrrolidinylpropyl) carboxamide;

3-((3-(4-fluorophenyl)(1H-indazol-5-yl))carbonylamino)-N-methyl propanamide;

3-((3-(4-fluorophenyl)(1H-indazol-5-yl))carbonylamino)-N,N-dimethyl propanamide;

3-((3-(4-fluorophenyl)(1H-indazol-5-yl))carbonylamino)-N-(2-methoxyethyl)propanamide; or a pharmaceutically acceptable salt thereof.

109. (Previously presented) A compound of claim 12, wherein the compound is:

phenyl-N-(3-phenyl(1H-indazol-5-yl))carboxamide;

N-(3-phenyl(1H-indazol-5-yl))-2-pyridylcarboxamide;

methyl 4-(N-(3-phenyl-1H-indazol-5-yl)carbamoyl)benzoate;

4-(N-(3-phenyl-1H-indazol-5-yl)carbamoyl)benzoic acid;

(2-hydroxyphenyl)-N-(3-phenyl(1H-indazol-5-yl)carboxamide);

N-(3-(phenyl-1H-indazole-5-yl))acetamide;
 (4-aminophenyl)-N-(3-phenyl(1H-indazol-5-yl))carboxamide;
 (3-aminophenyl)-N-(3-phenyl(1H-indazol-5-yl))carboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl)) (2-methylphenyl)carboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))(2-methoxyphenyl)carboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))(4-phenylphenyl)carboxamide;
 benzo(b)thiophen-2-yl-N-(3-(4-fluorophenyl)(1H-indazol-5-yl))carboxamide;
 methyl 4- {N-(3-(4-fluorophenyl)-1H-indazol-5-yl)carbamoyl} benzoate;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))-2-pyridylcarboxamide;
 4- {N-(3-(4-fluorophenyl)-1H-indazol-5-yl)carbamoyl} benzoic acid;
 cyclopropyl-N-(3-(4-fluorophenyl)(1H-indazol-5-yl))carboxamide;
 methyl 4- {N-(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-methylcarbamoyl} benzoate;
 4- {N-(3-fluorophenyl)(1H-indazol-5-yl))-N-methylcarbamoyl} benzoic acid;
 methyl 3- {N-((4-fluorophenyl)-1H-indazol-5-yl} carbamoyl} benzoate;
 3- {N-(3-(4-fluorophenyl)-1H-indazol-5-yl)carbamoyl} benzoic acid;
 N-(3-(4-fluorophenyl)-(1H-indazol-5-yl))(4-(N-methylcarbamoyl)phenyl)carboxamide;
 4- {N-(3-(4-fluorophenyl)-1H-indazol-5-yl)carbamoyl} benzamide;
 1-4- {N-(3-(4-methoxyphenyl)-1H-indazol-5-yl)carbamoyl} benzoic acid;
 4-(N-(3-(4-pyridyl)-1H-indazol-5-yl)carbamoyl)benzoic acid;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))benzamide;
 (3,4-bis(trifluoromethyl)phenyl)-N-(3-(4-fluorophenyl)(1H-indazol-5-yl))carboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))-2-furylcarboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))(3,4-dichlorophenyl)carboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))(2-hydroxyphenyl)carboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))-4-pyridylcarboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))-3-pyridylcarboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))-2-thienylcarboxamide;
 N-(3-(4-fluorophenyl)(1H-indazol-5-yl))morpholin-4-yl-carboxamide;
 [N-(((2R)-2-hydroxycyclohexyl)methyl) (3-(4-fluorophenyl) (1H-indazol-5-yl))carboxamide;] or a pharmaceutically acceptable salt thereof.

114. (Previously presented) A compound of claim 18, wherein the compound is:

3-(3-(4-fluorophenyl)-1H-indazol-5-yl)-1H-1,2,4-triazole;
5-(3-(4-fluorophenyl)(1H-indazole-5-yl))-3-methyl-4H-1,2,4-triazole;
5-{3-(4-fluorophenyl)(1H-indazole-5-yl)]-3-(methylethyl)-4H-1,2,4-triazole;
1-{5-(3-(4-fluorophenyl)-1H-indazole-5-yl)-4H-1,2,4-triazol-3-yl} propan-2-ol;
5-(3-(4-fluorophenyl)(1H-indazol-5-yl))-3-propyl-4H-1,2,4-triazole;
5-{3-(3-(methylethyl)phenyl)-1H-indazol-5-yl}-1H-1,2,4-triazole;
4-(5-(1H-1,2,4-triazol-5-yl)-1H-indazol-3-yl)phenol;
(4-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)dimethylamine;
{2-(4-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenoxy)ethyl} dimethylamine;
3-(5-(1H-1,2,4-triazol-5-yl)-1H-indazol-3-yl)furan;
1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-4-methoxybenzene;
5-(3-naphthyl-1H-indazol-5-yl)-1H-1,2,4-triazole;
3-(5-(1H-1,2,4-triazol-3-yl)-1H-indazol-3-yl)thiophene;
5-(3-(2-naphthyl)-1H-indazol-5-yl)-1H-1,2,4-triazole;
3-(5-(1H-1,2,4-triazol-3-yl)-1H-indazol-3-yl)phenylamine;
3-(3-(3,4-dichlorophenyl)-1H-indazol-5-yl)-1H-1,2,4-triazole;
3-(5-(1H-1,2,4-triazol-5-yl)-1H-indazol-3-yl)benzo(b)thiophene;
3-(3-(4-methylphenyl)-1H-indazol-5-yl)-1H-1,2,4-triazole;
N-(3-(5-(1H-1,2,4-triazol-3-yl)-1H-indazol-3-yl)phenyl)acetamide;
5-(3-(3-chlorophenyl)-1H-indazol-5-yl)-1H-1,2,4-triazole;
2-(5-(1H-1,2,4-triazol-5-yl)-1H-indazol-3-yl)benzo(b)furan;
1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-4-(methylsulfonyl)benzene;
1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-4-(methylsulfinyl)benzene;
5-(5-(1H-1,2,4-triazol-5-yl)-1H-indazol-3-yl)-2H-benzo(d)1,3-dioxolene;
4-(5-(1H-1,2,4-triazol-5-yl)-1H-indazol-3-yl)phenylamine;
5-{3-(4-(trifluoromethyl)phenyl)-1H-indazol-5-yl}-1H-1,2,4-triazole;
(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl)) phenyl) (methylsulfonyl)amine;
N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-methoxyacetamide;
N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-phenylacetamide;
N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-furylcarboxamide;
5-(3-(2-phenylethynyl)-1H-indazol-5-yl)-1H-1,2,4-triazole;

N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-3-pyridylcarboxamide;
 1-{5-{3-(4-fluorophenyl)1H-indazol-5-yl}-4H-1,2,4-Triazol-3-yl}ethan-1-ol;
 1-{5-(3-(4-fluorophenyl)-1H-indazol-5-yl)-4H-1,2,4-triazol-3-yl}propan-2-ol;
 {3-(3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenoxy)propyl}dimethylamine;
 {2-(3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenoxy)ethyl}dimethylamine;
 1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-3-(2-morpholin-4-yl-ethoxy)benzene;
 1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-3-(2-pyrrolidinylethoxy)benzene;
 1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-3-(2-piperidylethoxy)benzene;
 1-{2-(3-(5-(1H-1,2,4-triazol-5-yl)-1H-indazol-3-yl)phenoxy)ethyl}pyrrolidin-2-one;
 1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-3-(2-piperazinylethoxy)benzene;
 1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-3-(3-piperdylpropoxy)benzene;
 4-{2-(3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenoxy)ethyl}-1-acetylpiperazine;
 2-(3-(5-(1H-1,2,4-triazol-5-yl)-1H-indazol-3-yl)phenoxy)ethylamine;
 1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-3-(2-cyclohexylethoxy)benzene;
 1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-3-(2-azaperhyroepinylethoxy)benzene;
 N-(4-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-2-furyl caroxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-benzyl caroxamide;
 5-(3-(2-chlorophenyl)-1H-indazol-3-yl)-1H-1,2,4-triazole;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-(2,2-dimehtylpropyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-(cyclopropylmethyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-(3-pyridylmethyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-4-methyl piperazinyl ketone;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-((4-fluorophenyl)methyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-indan-2-ylcarboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-((1R)indanyl)carboxamide;

(3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-
 ((1S)indanyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-((1S,2R)-2-
 hydroxyindanyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-((2S,1R)-2-
 hydroxyindanyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-(1-methyl-1-
 phenylethyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-(tert-butyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-((1R)-1-
 phenylethyl)carboxamide;
 1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-3-(2-piperidylethoxy) benzene;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-isoindolin-2-yl ketone;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-(2-(dimethylamino)
 ethyl)carboxamide;
 1-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))-3-(2-piperidylethoxy) benzene;
 (3-(5-(1H-1,2,4-triazol-5-yl)(1H-indazol-3-yl))phenyl)-N-(1R)indanyl benzene;
 {5-(3-(4-fluorophenyl)-1H-indazol-5-yl)-4H-(1,2,4)-triazol-3-ylmethyl}-dimethyl-
 amine;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-3-piperidylpropanamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-hydroxypropanamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-
 (dimethylamino)acetamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)-1H-indazol-3-yl)phenyl)butanamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-phenoxypropanamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-3,3-dimethylbutanamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)cyclopropylcarboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)(6-chloro(3-
 pyridyl))carboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)cyclopentylcarboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)benzo(b)thiophen-2-
 carboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-pyridylcarboxamide;

N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-3-furylcarboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-hydroxy-2-phenylacetamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)isoxazol-5-ylcarboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)pentanamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-4-pyridylcarboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-cyclohexylacetamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-3-propanamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-(4-fluorophenyl)acetic acid;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)(2R)-2-hydroxy-2-phenylacetamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)(2S)-2-hydroxy-2-phenylacetamide;
 (2-{3-(3-(4-fluorophenyl)(1H-indazol-5-yl))(1H-1,2,4-triazol-5-yl)}ethyl)dimethylamine;
 diethyl({3-(3-(4-fluorophenyl)(1H-indazol-5-yl))(1H-1,2,4-triazol-5-yl)}methyl)amine;
 ({3-(3-(4-fluorophenyl)(1H-indazol-5-yl))(1H-1,2,4-triazol-5-yl)}methyl)methylamine;
 ({3-(3-(4-fluorophenyl)(1H-indazol-5-yl))(1H-1,2,4-triazol-5-yl)}ethyl)dimethylamine;
 (2R)-N-(3-(5-{5-((dimethylamino)methyl)(1H-1,2,4-triazol-3-yl)}(1H-indazol-3-yl))phenyl)-2-hydroxy-2-phenylacetamide;
 N-(3-(5-{5-((dimethylamino)methyl)(1H-1,2,4-triazol-3-yl)}(1H-indazol-3-yl))phenyl)-3,3-dimethylbutanamide;
 N-(3-(5-{5-((dimethylamino)methyl)(1H-1,2,4-triazol-3-yl)}(1H-indazol-3-yl))phenyl)-3-methylbutanamide;
 N-(3-(5-{5-((dimethylamino)methyl)(1H-1,2,4-triazol-3-yl)}(1H-indazol-3-yl))phenyl)-3-pyridylcarboxamide;
 (3-(5-{5-((dimethylamino)methyl)(1H-1,2,4-triazol-3-yl)}(1H-indazol-3-yl))phenyl)-N-((4-fluorophenyl)methyl)carboxamide;

(3-(5-{5-((dimethylamino)methyl)(1H-1,2,4-triazol-3-yl)}(1H-indazol-3-yl))phenyl)-
 N-((*tert*-butyl)methyl)carboxamide;
 ((1R)indanyl)(3-(5-{5-((dimethylamino)methyl)(1H-1,2,4-triazol-3-yl)}(1H-indazol-
 3-yl))phenyl)carboxamide;
 ({3-(3-(4-methoxyphenyl)(1H-indazol-5-yl))(1H-1,2,4-triazol-5-
 yl)}methyl)dimethylamine;
 {(3-(3-(2H-benzo(d)1,3-dioxolen-5-yl))(1H-indazol-5-yl))(1H-1,2,4-triazol-5-
 yl)}methyl}dimethylamine;
 (3-(5-(3-((dimethylamino)methyl)(1H-1,2,4-triazol-5-yl))(1H-indazol-3-yl))phenyl)-
 N-(2-piperidylethyl)carboxamide;
 ((5-(3-benzo(D)furan-2-yl(1H-indazol-5-yl))(1H-1,2,4-triazol-3-
 yl))methyl)dimethylamine;
 (3-(5-(3-((dimethylamino)methyl)(1H-1,2,4-triazol-5-yl))(1H-indazol-3-yl))phenyl)-
 N-benzamide;
 (3-(5-(3-((dimethylamino)methyl)(1H-1,2,4-triazol-5-yl))(1H-indazol-3-yl))phenyl)-
 N-(4-fluorophenyl)carboxamide-2HCl;
 (3-(5-(3-((dimethylamino)methyl)(1H-1,2,4-triazol-5-yl))(1H-indazol-3-yl))phenyl)-
 N-indan-2-yl-carboxamide;
 (3-(5-(3-((dimethylamino)methyl)(1H-1,2,4-triazol-5-yl))(1H-indazol-3-yl))phenyl)-
 N-cyclopropylcarboxamide;
 (3-(5-(3-((dimethylamino)methyl)(1H-1,2,4-triazol-5-yl))(1H-indazol-3-yl))phenyl)-
 N-cyclobutylcarboxamide-2HCl;
 1-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl)-3-(2-methoxyethoxy)benzene;
 1-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl)-3-(3-pyridylmethoxy)benzene;
 3-(5-(1H-1,2,4-triazol-3-yl)-1H-indazol-3-yl)benzoic acid;
 3-(5-(1H-1,2,4-triazol-3-yl)-1H-indazol-3-yl)benzoic acid N-(4-(5-(1H-1,2,4-triazol-
 3-yl)(1H-indazol-3-yl))phenyl)-2-(3-pyridyl)acetamide;
 N-(4-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-phenylacetamide;
 N-(4-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-methoxyacetamide;
 N-(4-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-
 (dimethylamino)acetamide;
 (4-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)(methylsulfonyl)amine;

(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(2-methoxyethyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-benzamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(2-phenethyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(2-piperidylethyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(2-morpholin-4-ylethyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-cyclohexylcarboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-cyclopentylcarboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(4-fluorophenyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-((1R,2R)-2-phenylcyclopropyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-cyclopropylcarboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(3-pyridyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(5,6,7,8-tetrahydronaphthyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(1-benzyl(4-piperidyl))carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(1-benzylpyrrolidin-3-yl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(methylethyl)carboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-cyclobutylcarboxamide;
 (3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-N-(4-pyridyl)carboxamide;
 6-(5-(1H-1,2,4-triazol-3-yl)-1H-indazol-3-yl)-2H,3h-benzo(e)1,4-dioxin;
 6-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))-2-methoxynaphthalene;
 3-(3-(3-quinoyl)-1H-indazol-5-yl)-1H-1,2,4-triazole;
 5-(5-(1H-1,2,4-triazol-3-yl)-1H-indazol-3-yl)-2,3-dihydrobenzo(b)furan;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)benzamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)(2,4-dichlorophenyl)carboxamide;

N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)(4-methoxyphenyl)carboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)(4-methylphenyl)carboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)(4-chlorophenyl)carboxamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-methylpropanamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-3-methylbutanamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-morpholin-4-yl-acetamide;
 N-(3-(5-(1H-1,2,4-triazol-3-yl)(1H-indazol-3-yl))phenyl)-2-(4-methylpiperazinyl)acetamide;
 2-methoxy-6-{5-(5-(pyrrolidinylmethyl)(1H-1,2,4-triazol-3-yl))(1H-indazol-3-yl)}naphthalene;
 N-phenyl(3-{5-(5-(pyrrolidinylmethyl)(1H-1,2,4-triazol-3-yl))(1H-indazol-3-yl)})carboxamide;
 6-{5-(5-(pyrrolidinylmethyl)-1H-1,2,4-triazol-3-yl)-1H-indazol-3-yl}-2H,3H-benzo(e)1,4-dioxin; or a pharmaceutically acceptable salt thereof.

115. (Previously presented) A compound of claim 19, wherein the compound is:

5-(3-(4-fluorophenyl)-1H-indazol-5-yl)-2H-1,2,3,4-tetrazole;
 1-(5-(2H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))-2-methoxybenzene;
 5-(3-(3-pyridyl)-1H-indazol-5-yl)-2H-1,2,3,4-tetrazole;
 2-(5-(2H-1,2,3,4-tetrazol-5-yl)-1H-indazol-3-yl)thiophene;
 5-{3-(4-(methylethyl)phenyl)-1H-indazol-5-yl}-2H-1,2,3,4-tetrazole;
 2-(5-(2H-1,2,3,4-tetrazol-5-yl)-1H-indazol-3-yl)furan;
 3-(5-(2H-1,2,3,4-tetrazol-5-yl)-1H-indazol-3-yl)phenylamine;
 5-(5-(1H-1,2,3,4-tetraazol-5-yl)-1H-indazol-3-yl)-2H-benzo(d)1,3-dioxolene;
 3-(5-(2H-1,2,3,4-tetrazol-5-yl)-1H-indazol-3-yl)thiophene;
 5-(3-(2-naphthyl)-1H-indazol-5-yl)-1H-1,2,3,4-tetrazole;
 1-(5-(1H-1,2,3,4-tetraazol-5-yl)(1H-indazol-3-yl))-4-methoxybenzene;
 1-(5-(1H-1,2,3,4-tetraazol-5-yl)(1H-indazol-3-yl))-4-(2-methylpropoxy)benzene;
 5-(3-(4-chlorophenyl)-1H-indazol-5-yl)-2H-1,2,3,4-tetrazole;

1-(5-(2H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))-3-methoxybenzene;
 5-(3-(4-pyridyl)-1H-indazol-5-yl)-2H-1,2,3,4-tetrazole;
 2-(5-(2H-1,2,3,4-tetraazol-5-yl)-1H-indazol-3-yl)benzo(b)furan;
 2-(5-(2H-1,2,3,4-tetrazol-5-yl)-1H-indazol-3-yl)phenol;
 3-(5-(2H-1,2,3,4-tetrazol-5-yl)-1H-indazol-3-yl)phenol;
 5-(3-(2-phenylethynyl)-1H-indazol-5-yl)-1H-1,2,3,4-tetrazole;
 5-(3-(2-phenylethyl)-1H-indazol-5-yl)-2H-1,2,3,4-tetrazole;
 N-(3-(5-(2H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))phenyl)-2-methoxyacetamide;
 2-(5-(1H-1,2,3,4-tetraazol-5-yl)-1H-indazol-3-yl)benzo(b)thiophene;
 1-(5-(1H-1,2,3,4-tetraazol-5-yl)(1H-indazol-3-yl))-4-(2-morpholin-4-
 ylethoxy)benzene;
 N-(3-(5-2H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))phenyl)-2-phenoxypropanamide;
 N-(3-(5-(2H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))phenyl)-3-
 piperidylpropanamide;
 N-(3-(5-(2H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))phenyl)-2-furylcarboxamide;
 1-(5-(1H-1,2,3,4-tetraazol-5-yl)(1H-indazol-3-yl))-3-(2-morpholin-4-
 ylethoxy)benzene;
 4-(5-(2H-1,2,3,4-tetraazol-5-yl)(1H-indazol-3-yl))-1,2-dimethoxybenzene;
 N-(3-(5-(2H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))phenyl)-3-
 methoxypropanamide;
 N-(3-(5-(2H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))phenyl)-3-pyridylcarboxamide;
 {3-(4-(5-(1H-1,2,3,4-tetrazo-5-yl)(1H-indazol-3-yl))phenoxy)propyl}
 dimethylamine;
 {3-(3-(5-(1H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))phenoxy)propyl}
 dimethylamine;
 {2-(3-(5-(1H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))phenoxy)ethyl}dimethylamine;
 N-(3-(5-(2H-1,2,3,4-tetrazol-5-yl)(1H-indazol-3-yl))phenyl)-(2S)-2-
 hydroxypropanamide;
 N-(4-(5-(2H-1,2,3,4-tetrazo-5-yl)(1H-indazol-3-yl))phenyl)-3-pyridylcarboxamide;
 or a pharmaceutically acceptable salt thereof.

116. (Previously presented)
 compound is:

A compound of claim 20, wherein the

3-(4-fluorophenyl)-5-imidazol-2-yl-1H-indazole, or a pharmaceutically acceptable salt thereof.

117. (Previously presented) A compound, wherein the compound is:
3-phenyl-5-(phenylmethoxy)-1H-indazole;
(3-(4-fluorophenyl)(1H-indazol-5-yl))(phenylsulfonyl)amine;
3-(4-fluorophenyl)-1H-indazole-5-carboxylate;
(3-(4-fluorophenyl)(1H-indazol-5-yl))-N-(phenylmethoxy)carboxamide;
3-(4-fluorophenyl)-1H-indazole-5-carbohydroxamic acid;
N-((tert-butoxy)carbonylamino) (3-(4-fluorophenyl) (1H-indazol-5-yl))carboxamide;
N-amino(3-(4-fluorophenyl)(1 H-indazol-5-yl))carboxamide;
methyl-3-benzo(B)thiophen-2-yl-1H-indazole-5-carboxylate;
3-benzo(B)thiophen-2-yl-1H-indazole-5-carboxylic acid;
or a pharmaceutically acceptable salt thereof.